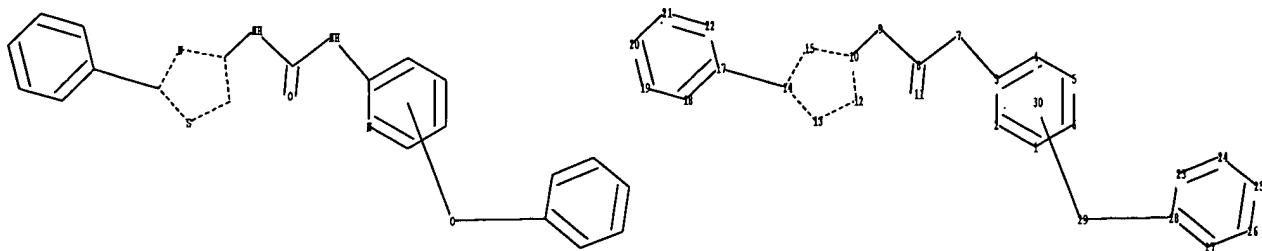


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chain nodes :

7 8 9 11 29

ring nodes :

1 2 3 4 5 6 10 12 13 14 15 17 18 19 20 21 22 23 24 25 26 27
28

chain bonds :

3-7 7-8 8-9 8-11 9-10 14-17 28-29

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 10-12 10-15 12-13 13-14 14-15 17-18 17-22
18-19 19-20 20-21 21-22 23-28 23-24 24-25 25-26 26-27 27-28

exact/norm bonds :

3-7 7-8 8-9 8-11 9-10 10-12 10-15 12-13 13-14 14-15 28-29

exact bonds :

14-17

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 17-18 17-22 18-19 19-20 20-21 21-22 23-28
23-24 24-25 25-26 26-27 27-28

isolated ring systems :

containing 1 : 10 : 17 : 23 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:Atom
11:CLASS 12:Atom 13:Atom 14:Atom 15:Atom 17:Atom 18:Atom 19:Atom 20:Atom
21:CLASS 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:CLASS
30:CLASS

L2 STRUCTURE UPLOADED

=> d

L2 HAS NO ANSWERS

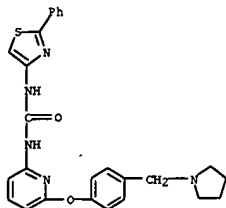
L2 STR

*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***

Structure attributes must be viewed using STN Express query preparation.

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L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 479365-32-1 REGISTRY
ED Entered STN: 08 Jan 2003
CN Urea, N-(2-phenyl-4-thiazolyl)-N'-[6-[4-(1-pyrrolidinylmethyl)phenoxy]-2-pyridinyl]- (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C26 H25 N5 O2 S
SR CA
LC STN Files: CA, CAPLUS, TOXCENTER, USPAT2, USPATFULL



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1 REFERENCES IN FILE CA (1907 TO DATE)
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

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```
=> d ibib abs hitstr tot
'IBIB' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
'ABS' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
'HITSTR' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'
```

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

```
REG      - RN
SAM      - Index Name, MF, and structure - no RN
FIDE     - All substance data, except sequence data
IDE      - FIDE, but only 50 names
SQIDE    - IDE, plus sequence data
SQIDE3   - Same as SQIDE, but 3-letter amino acid codes are used
SQD      - Protein sequence data, includes RN
SQD3     - Same as SQD, but 3-letter amino acid codes are used
SQN      - Protein sequence name information, includes RN

CALC     - Table of calculated properties
EPROP    - Table of experimental properties
PROP     - EPROP and CALC
```

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

```
ABS  -- Abstract
APPS -- Application and Priority Information
BIB  -- CA Accession Number, plus Bibliographic Data
CAN  -- CA Accession Number
CBIB -- CA Accession Number, plus Bibliographic Data (compressed)
IND  -- Index Data
IPC  -- International Patent Classification
PATS -- PI, SO
STD  -- BIB, IPC, and NCL
```

```
IABS -- ABS, indented, with text labels
IBIB -- BIB, indented, with text labels
ISTD -- STD format, indented
```

```
OBIB ----- AN, plus Bibliographic Data (original)
OIBIB ----- OBIB, indented with text labels
```

```
SBIB ----- BIB, no citations
SIBIB ----- IBIB, no citations
```

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help

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messages:

HELP DFIELDS -- To see a complete list of individual display fields.
HELP FORMATS -- To see detailed descriptions of the predefined formats.
ENTER DISPLAY FORMAT (IDE):ide

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L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2006 ACS on STN
RN 478365-32-1 REGISTRY
ED Entered STN: 08 Jan 2003
CN Urea, N-(2-phenyl-4-thiazolyl)-N'-[6-[4-(1-pyrrolidinylmethyl)phenoxy]-2-pyridinyl]- (9CI) (CA INDEX NAME)
FS 3D CONCORD
MF C26 H25 N5 O2 S
SR CA
LC STN Files: CA, CAPLUS, TOXCENTER, USPAT2, USPATFULL

